Welcome to STN International! Enter x:x

LOGINID:SSPTAVXR1614

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * *	* *	* *	* *	* Welcome to STN International * * * * * * * * * *
NEWS	1			Web Page for STN Seminar Schedule - N. America
NEWS	2	AUG	10	Time limit for inactive STN sessions doubles to 40
				minutes
NEWS	3	AUG	18	COMPENDEX indexing changed for the Corporate Source
NEWS		3.110	0.4	(CS) field ENCOMPLIT/ENCOMPLIT2 reloaded and enhanced
NEWS			24	
MEMP	,	AUG	24	U.S. patents
NEWS	6	SEP	09	50 Millionth Unique Chemical Substance Recorded in
				CAS REGISTRY
NEWS	7	SEP	11	WPIDS, WPINDEX, and WPIX now include Japanese FTERM
				thesaurus
NEWS	8	OCT	21	Derwent World Patents Index Coverage of Indian and
NIB110		OCT	0.1	Taiwanese Content Expanded
NEWS	9	OCI	21	Derwent World Patents Index enhanced with human translated claims for Chinese Applications and
				Utility Models
NEWS	10	OCT	27	Free display of legal status information in CA/CAplus,
			_	USPATFULL, and USPAT2 in the month of November.
NEWS	EXP	RESS		26 09 CURRENT WINDOWS VERSION IS V8.4,
			AND	CURRENT DISCOVER FILE IS DATED 06 APRIL 2009.
NEWS	HOII	RS	STI	N Operating Hours Plus Help Desk Availability
NEWS				lcome Banner and News Items
				ed by the item number or name to see news on that
speci	fic	topi	٥.	
3.11			n men r	in and the terminal and the committee of the committee of
				is subject to the provisions of the STN customer agreement limits use to scientific research. Use
				velopment or design, implementation of commercial
				e of CAS and STN data in the building of commercial
, , , ,	- 4			

products is prohibited and may result in loss of user privileges

FILE 'HOME' ENTERED AT 14:23:13 ON 18 NOV 2009

and other penalties.

=> file registry SINCE FILE TOTAL ENTRY SESSION 0.22 0.22 COST IN U.S. DOLLARS FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 14:23:25 ON 18 NOV 2009

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STRUCTURE FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8 DICTIONARY FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 26, 2009.

Please note that search-term pricing does apply when conducting ${\tt SmartSELECT}$ searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

=> file hcaplus

 COST IN U.S. DOLLARS
 SINCE FILE
 TOTAL

 FULL ESTIMATED COST
 0.48
 0.70

FILE 'HCAPLUS' ENTERED AT 14:24:18 ON 18 NOV 2009
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FILE COVERS 1907 - 18 Nov 2009 VOL 151 ISS 21
FILE LAST UPDATED: 17 Nov 2009 (20091117/ED)
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Aug 2009
USPIO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Aug 2009

 ${\tt HCAplus}$ now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2009.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

During November, try the new LSUS format of legal status information in the ${\rm CA/CAplus}$ family databases for free! Complete details on the

number of free displays and other databases participating in this offer appear in NEWS 10.

=> e us 2006566384

NUMBER OF TERMS TO DISPLAY IS OUT OF RANGE

The total number of terms displayed in a single EXPAND command must be in the range 5-25.

```
=> e us 2006566384/apps
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EI	1	UA99-95363/PRN
E2	1	UA99-98052573/PRN
E3	0>	US 2006566384/AP
E 4	0	US 2006566384/PRN
E5	1	US00-10056/AP
E6	1	US00-10070/AP
E7	1	US00-10178/AP
E8	1	US00-10197/AP
E9	3	US00-10220/AP
E10	1	US00-102457/AP
E11	1	US00-10268/AP
E12	1	US00-103430/AP

=> e us 2006-566384/apps

-/ e us	2000-30030	1/appo
E1	2	US2006-566383/AP
E2	1	US2006-566383/PRN
E3	1>	US2006-566384/AP
E4	0	US2006-566384/PRN
E5	2	US2006-566385/AP
E6	1	US2006-566387/AP
E7	1	US2006-566388/AP
E8	1	US2006-566389/AP
E9	1	US2006-566390/AP
E10	1	US2006-566390/PRN
E11	2	US2006-566392/AP
E12	1	US2006-566393/AP

=> s us 2006-566384/apps

- 1 US2006-566384/AP
- 0 US2006-566384/PRN 1 US 2006-566384/APPS

(US2006-566384/AP, PRN)

=> sel rn

E1 THROUGH E6 ASSIGNED

=> fi	.le	regis	stry
COST	IN	U.S.	DOLLARS

COST	IN U.S.	DOLLARS	SINCE FILE	TOTAL
			ENTRY	SESSION
FULL	ESTIMATE	D COST	5.80	6.50

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STRUCTURE FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8 DICTIONARY FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 26, 2009.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

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http://www.cas.org/support/stngen/stndoc/properties.html

```
=> s e1-e6
             1 1044645-80-8/BI
                 (1044645-80-8/RN)
             1 195875-87-7/BI
                 (195875-87-7/RN)
             1 50-36-2/BI
                 (50-36-2/RN)
             1 54-11-5/BI
                 (54-11-5/RN)
             1 835651-44-0/BI
                 (835651-44-0/RN)
             1 87-69-4/BI
                 (87-69-4/RN)
T.2
             6 (1044645-80-8/BI OR 195875-87-7/BI OR 50-36-2/BI OR 54-11-5/BI
               OR 835651-44-0/BI OR 87-69-4/BI)
```

=> d hitstr

- RN

REG

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data

DE - FIDE, but only 50 names
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence name information, includes RN
EPROP - Table of experimental properties
PPROP - Table of predicted properties
PPROP - Table of predicted properties

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

```
ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
```

IPC -- International Patent Classification PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original) OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

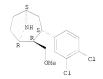
The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats. ENTER DISPLAY FORMAT (IDE):ide THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) /N:y

- ANSWER 1 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN
- 1044645-80-8 REGISTRY RN
- Entered STN: 29 Aug 2008
- CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,
- (1R, 2R, 3S, 5S) -rel- (CA INDEX NAME) FS STEREOSEARCH
- ME C15 H19 C12 N O
- SR CA
- LC. STN Files: CA, CAPLUS, TOXCENTER

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE) 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 2 'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY' The following are valid formats: Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number) - RN SAM - Index Name, MF, and structure - no RN - All substance data, except sequence data - FIDE, but only 50 names SQIDE - IDE, plus sequence data SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used SOD - Protein sequence data, includes RN SOD3 - Same as SQD, but 3-letter amino acid codes are used SON - Protein sequence name information, includes RN EPROP - Table of experimental properties PPROP - Table of predicted properties PROP - EPROP, ETAG, PPROP and SPEC Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are: ABS -- Abstract APPS -- Application and Priority Information BIB -- CA Accession Number, plus Bibliographic Data CAN -- CA Accession Number CBIB -- CA Accession Number, plus Bibliographic Data (compressed) IND -- Index Data IPC -- International Patent Classification PATS -- PI, SO STD -- BIB, IPC, and NCL IABS -- ABS, indented, with text labels IBIB -- BIB, indented, with text labels ISTD -- STD format, indented OBIB ----- AN, plus Bibliographic Data (original) OIBIB ----- OBIB, indented with text labels SBIB ----- BIB, no citations SIBIB ----- IBIB, no citations The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available. The MAX format is the same as ALL. The IALL format is the same as ALL with BIB ABS and IND indented, with text labels. For additional information, please consult the following help

messages:

HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats. ENTER DISPLAY FORMAT (IDE):ide THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) /N:y

ANSWER 2 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN RN 835651-44-0 REGISTRY

ED Entered STN: 22 Feb 2005

8-Azabicyclo[3,2,1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,
(1R,2R,3S,5S)-, (2R,3R)-2,3-dihydroxybutanedioate (1:1) (CA INDEX NAME)

STERSOSEARCH

C15 H19 C12 N O . C4 H6 O6

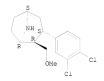
SR CA

CSTN Files: CA, CAPLUS, EMBASE, TOXCENTER, USPATFULL

CM 1

CRN 195875-87-7 CMF C15 H19 C12 N O

Absolute stereochemistry.



CM

CRN 87-69-4 CMF C4 H6 O6

Absolute stereochemistry.

2 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 3

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

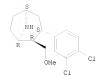
SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names

SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

```
SQD
       - Protein sequence data, includes RN
SQD3
       - Same as SQD, but 3-letter amino acid codes are used
SON
       - Protein sequence name information, includes RN
EPROP - Table of experimental properties
PPROP - Table of predicted properties
PROP
      - EPROP, ETAG, PPROP and SPEC
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ISTD -- STD format, indented
OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N:v
     ANSWER 3 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN
     195875-87-7 REGISTRY
RN
     Entered STN: 23 Oct 1997
ED
     8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,
     (1R, 2R, 3S, 5S) - (CA INDEX NAME)
OTHER CA INDEX NAMES:
    8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,
     [1R-(2-endo, 3-exo)]-
FS
    STEREOSEARCH
MF
    C15 H19 C12 N O
CT
    COM
SR
LC
     STN Files: CA, CAPLUS, CASREACT, EMBASE, TOXCENTER, USPATZ, USPATFULL
Absolute stereochemistry.
```



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

6 REFERENCES IN FILE CA (1907 TO DATE) 6 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 4 'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RM

SAM - Index Name, MF, and structure - no RN

FIDE - All substance data, except sequence data

- FIDE, but only 50 names IDE

SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

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PPROP - Table of predicted properties

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CBIB -- CA Accession Number, plus Bibliographic Data (compressed) IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

```
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
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The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) /N:y
     ANSWER 4 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN
RN
     87-69-4 REGISTRY
ED
     Entered STN: 16 Nov 1984
    Butanedioic acid, 2,3-dihydroxy- (2R,3R)- (CA INDEX NAME)
OTHER CA INDEX NAMES:
    Butanedioic acid, 2,3-dihydroxy- [R-(R*,R*)]-
     Tartaric acid, L-(+)- (8CI)
OTHER NAMES:
CN
    (+)-(2R,3R)-Tartaric acid
CN
     (+)-(R,R)-Tartaric acid
CN
    (+)-L-Tartaric acid
CN
     (+)-Tartaric acid
CN
    (2R,3R)-(+)-Tartaric acid
CN
    (2R,3R)-2,3-Dihydroxysuccinic acid
CN
    (2R,3R)-Tartaric acid
CN
    (R,R)-(+)-Tartaric acid
CN
    (R,R)-Tartaric acid
CN
    1,2-Dihydroxyethane-1,2-dicarboxylic acid
CN
    2,3-Dihydroxybutanedioic acid
CN
    d-α,β-Dihydroxysuccinic acid
CN
    d-Tartaric acid
CN
    Dextrotartaric acid
CN
    Dihydroxysuccinic acid
CN
    E 334
CN
     L-(+)-Tartaric acid
CN
    L-Tartaric acid
CN
    Natural tartaric acid
    NSC 62778
CN
CN
     Tartaric acid
CN
     Threaric acid
DR
     1039646-76-8, 8014-54-8, 8059-77-6, 1336-18-1
MF
     C4 H6 O6
     COM
LC
     STN Files:
                ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS,
       BIOTECHNO, CA, CABA, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMINFORMRX,
       CHEMLIST, CIN, CSCHEM, DDFU, DETHERM*, DRUGU, EMBASE, ENCOMPLIT,
       ENCOMPLIT2, ENCOMPPAT, ENCOMPPAT2, GMELIN*, IFICDB, IFIPAT, IFIUDB, IPA,
       MRCK*, MSDS-OHS, NAPRALERT, PIRA, PROMT, PS, RTECS*, SPECINFO,
       SYNTHLINE, TOXCENTER, TULSA, USAN, USPAT2, USPATFULL, USPATOLD
         (*File contains numerically searchable property data)
     Other Sources: DSL**, EINECS**, TSCA**
```

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

25806 REFERENCES IN FILE CA (1907 TO DATE)

2072 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 25890 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 5

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

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- FIDE, but only 50 names

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SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

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PPROP - Table of predicted properties

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APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed) IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

```
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) /N:y
    ANSWER 5 OF 6 REGISTRY COPYRIGHT 2009 ACS on STN
RN
    54-11-5 REGISTRY
ED
    Entered STN: 16 Nov 1984
    Pyridine, 3-((2S)-1-methyl-2-pyrrolidinyl)- (CA INDEX NAME)
OTHER CA INDEX NAMES:
   Nicotine (8CI)
    Pyridine, 3-(1-methyl-2-pyrrolidinyl)-, (S)-
OTHER NAMES:
CN
    (-)-(S)-Nicotine
CN
    (-)-β-Pyridyl-α-N-methylpyrrolidine
CN
    (-)-3-(1-Methyl-2-pyrrolidyl)pyridine
CN
    (-)-Nicotine
CN
    (S)-(-)-Nicotine
CN
    (S)-3-(1-Methyl-2-pyrrolidinyl)pyridine
CN
    (S)-Nicotine
CN
    3-[(2S)-1-Methyl-2-pyrrolidinyl]pyridine
CN
    Exodus
CN Flux Maag
CN
    Habitrol
CN
   L-Nicotine
CN
    1-Nicotine
CN
   Nicabate
CN
   Nicoderm
CN
   Nicolan
CN
    Niconil
CN
    Nicopatch
CN
    Nicorette
CN
    Nicotell TTS
CN
    Nicotin
CN
    Nicotinell
CN
    Nicotrol
CN
    Nikofrenon
CN
    Niquitin
CN
    NSC 5065
CN
    Tabazur
CN
    XL All Insecticide
FS
    STEREOSEARCH
DR
    13890-81-8, 13890-82-9, 6912-85-2, 551-13-3, 16760-37-5
ME
    C10 H14 N2
    COM
    STN Files:
                ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*,
      BIOSIS, BIOTECHNO, CA, CABA, CAPLUS, CASREACT, CBNB, CHEMCATS,
```

CHEMINFORMEX, CHEMLIST, CHEMSAFE, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IMSCOSEARCH, IMSDBUGNEWS, IMSPATENTS, IMSRESEARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, NAFRALERT, PHAR, PIRA, PROMT, RTECS*, SPECINFO, TOXCENTER, ULIDAT, USAN, USPATZ), USPATOLU, VETU

(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**

(**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry. Rotation (-).

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

22661 REFERENCES IN FILE CA (1907 TO DATE)
395 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
22714 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d hitstr 6

'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN

FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names

SOIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

EPROP - Table of experimental properties

PPROP - Table of predicted properties

PROP - EPROP, ETAG, PPROP and SPEC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

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STD -- BIB, IPC, and NCL
IABS -- ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented
OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.
For additional information, please consult the following help
messages:
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide
THE ESTIMATED COST FOR THIS REQUEST IS 2.05 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) /N:v
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RN
    50-36-2 REGISTRY
ED
    Entered STN: 16 Nov 1984
    8-Azabicyclo[3.2.1]octane-2-carboxylic acid, 3-(benzoyloxy)-8-methyl-,
    methyl ester, (1R, 2R, 3S, 5S) - (CA INDEX NAME)
OTHER CA INDEX NAMES:
    1αH,5αH-Tropane-2β-carboxylic acid, 3β-hydroxy-,
     methyl ester, benzoate (ester) (8CI)
CN
     8-Azabicyclo[3.2.1]octane-2-carboxylic acid, 3-(benzoyloxy)-8-methyl-,
    methyl ester, [1R-(exo,exo)]-
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                    EINECS**
     Other Sources:
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(**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry. Rotation (-).

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

15922 REFERENCES IN FILE CA (1907 TO DATE) 333 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 15948 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d his

L2

(FILE 'HOME' ENTERED AT 14:23:13 ON 18 NOV 2009)

FILE 'REGISTRY' ENTERED AT 14:23:25 ON 18 NOV 2009

FILE 'HCAPLUS' ENTERED AT 14:24:18 ON 18 NOV 2009

E US 2006566384/APPS E US 2006-566384/APPS

L1 1 S US 2006-566384/APPS SEL RN

6 S E1-E6

FILE 'REGISTRY' ENTERED AT 14:25:37 ON 18 NOV 2009

=> s 835651-44-0/rn

L3 1 835651-44-0/RN

=> file hcaplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 22.38 28.88

FILE 'HCAPLUS' ENTERED AT 14:37:58 ON 18 NOV 2009 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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FILE COVERS 1907 - 18 Nov 2009 VOL 151 ISS 21 FILE LAST UPDATED: 17 Nov 2009 (20091117/ED) REVISED CLASS FIELDS (/NCL) LAST RELOADED: Aug 2009 USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Aug 2009

HCAplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2009.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

During November, try the new LSUS format of legal status information in the CA/CAplus family databases for free! Complete details on the number of free displays and other databases participating in this offer appear in NEWS 10.

=> s 13 L4 2 L3

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THE ESTIMATED COST FOR THIS REQUEST IS 7.78 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N:y

L4 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2009 ACS on STN ACCESSION NUMBER: 2008:1070913 HCAPLUS

DOCUMENT NUMBER: 149:315512

TITLE: Hydrochloride salt of an azabicyclo[3.2.1]octane

derivative
INVENTOR(S): Grinter, Trevor John; Moldt, Peter; Watjen, Frank

PATENT ASSIGNEE(S): Glaxo Group Limited, UK; Neurosearch A/S

SOURCE: PCT Int. Appl., 50pp.
CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English FAMILY ACC. NUM. COUNT: 1

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OTHER SOURCE(S): CASREACT 149:315512

IT 835651-44-0

RL: RCT (Reactant); RACT (Reactant or reagent)
(hydrochloride salt of an azabicyclo[3.2.1]octane derivative)

RN 835651-44-0 HCAPLUS

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CRN 195875-87-7 CMF C15 H19 C12 N O

Absolute stereochemistry.

CM :

CRN 87-69-4 CMF C4 H6 O6

Absolute stereochemistry.

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REFERENCE COUNT:

THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2009 ACS on STN ACCESSION NUMBER: 2005:120742 HCAPLUS

DOCUMENT NUMBER: 142:183325

TITLE: Preparation of

2-methoxymethyl-3-(3,4-dichlorophenyl)-8azabicyclo[3.2.1]octane tartrate salts

INVENTOR(S): Frostrup, Brian; Waetjen, Frank; Jensen, Klaus Snej
PATENT ASSIGNEE(S): Neurosearch A/S, Den.

SOURCE: PCT Int. Appl., 19 pp.

DOCUMENT TYPE: Patent

LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1

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IT 835651-44-0P

RL: PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of methoxymethyldichlorophenyl azabicyclooctane tartrate salts)
RN 835651-44-0 HCAPLUS

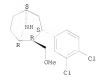
CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R,2R,3S,5S)-, (2R,3R)-2,3-dihydroxybutanedioate (1:1) (CA INDEX NAME)

CM 1

CRN 195875-87-7

CMF C15 H19 C12 N O

Absolute stereochemistry.



CM 2

CRN 87-69-4 CMF C4 H6 O6

Absolute stereochemistry.

OH OH

OS.CITING REF COUNT:

- 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD (4 CITINGS)
- REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> file registry

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 10.63 39.51

FULL ESTIMATED COST

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8 DICTIONARY FILE UPDATES: 16 NOV 2009 HIGHEST RN 1192511-54-8

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TSCA INFORMATION NOW CURRENT THROUGH June 26, 2009.

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http://www.cas.org/support/stngen/stndoc/properties.html

=> s 1044645-80-8/rn or 195875-87-7/rn

- 1 1044645-80-8/RN
- 1 195875-87-7/RN
- L5 2 1044645-80-8/RN OR 195875-87-7/RN

=> file hcaplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST Session 0.48 39.99

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FILE COVERS 1907 - 18 Nov 2009 VOL 151 ISS 21 FILE LAST UPDATED: 17 Nov 2009 (20091117/ED) REVISED CLASS FIELDS (/NCL) LAST RELOADED: Aug 2009 USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Aug 2009

HCAplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2009.

CAS Information Use Policies apply and are available at:

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This file contains CAS Registry Numbers for easy and accurate substance identification.

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=> s 15

L7 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2009 ACS on STN ACCESSION NUMBER: 2005:120742 HCAPLUS

DOCUMENT NUMBER: 142:183325 TITLE: Preparation of

2-methoxymethyl-3-(3,4-dichlorophenyl)-8azabicyclo[3.2.1]octane tartrate salts

Frostrup, Brian; Waetjen, Frank; Jensen, Klaus Snej INVENTOR(S):

Neurosearch A/S, Den. PATENT ASSIGNEE(S): SOURCE: PCT Int. Appl., 19 pp.

CODEN: PIXXD2 DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1 PATENT INFORMATION:

PATENT NO.

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THE ESTIMATED COST FOR THIS REQUEST IS 19.45 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y) /N:y

L12 ANSWER 1 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2008:1070913 HCAPLUS

DOCUMENT NUMBER: 149:315512

TITLE: Hydrochloride salt of an azabicyclo[3.2.1]octane derivative

INVENTOR(S): Grinter, Trevor John; Moldt, Peter; Watjen, Frank PATENT ASSIGNEE(S): Glaxo Group Limited, UK; Neurosearch A/S

SOURCE: PCT Int. Appl., 50pp.

CODEN: PIXXD2 DOCUMENT TYPE: Patent

LANGUAGE: English FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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OTHER SOURCE(S): CASREACT 149:315512

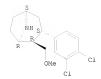
IT 195875-87-7P RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(hydrochloride salt of an azabicyclo[3.2.1]octane derivative)

195875-87-7 HCAPLUS RN

8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, CN (1R, 2R, 3S, 5S) - (CA INDEX NAME)

Absolute stereochemistry.



REFERENCE COUNT:

6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 2 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2007:478903 HCAPLUS

DOCUMENT NUMBER: 146:468550

TITLE: Treatment of diabetes with tropanes

Patent

INVENTOR(S): Dugi, Klaus; Berger, Frank; Raschig, Andreas; Reess,
Juergen; Salin, Laurence

PATENT ASSIGNEE(S): Boehringer Ingelheim Pharma G.m.b.H. & Co. K.-G.,

Germany SOURCE: Eur. Pat. Appl., 19pp.

CODEN: EPXXDW

DOCUMENT TYPE:

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1 PATENT INFORMATION:

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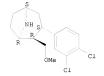
OTHER SOURCE(S): MARPAT 146:468550 TT 195875-87-7

RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses) (treatment of diabetes with tropanes)

195875-87-7 HCAPLUS RN

CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R, 2R, 3S, 5S) - (CA INDEX NAME)

Absolute stereochemistry.



OS.CITING REF COUNT: THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD 2 (2 CITINGS)

REFERENCE COUNT: 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 3 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN 2003:971889 HCAPLUS

ACCESSION NUMBER:

DOCUMENT NUMBER: 140:13076

TITLE: Triple monoamine reuptake inhibitors for the treatment

of chronic pain

INVENTOR(S): Scheel-Krueger, Jorgen; Blackburn-Munro, Gordon John

PATENT ASSIGNEE(S): Neurosearch A/S, Den. SOURCE: PCT Int. Appl., 20 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Pat.ent. English

LANGUAGE: FAMILY ACC. NUM. COUNT: 1

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PA:	TENT :	NO.			KIN	D	DATE		APPLICATION NO.									
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		GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	KZ,	LC,	LK,	LR,	
		LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NI,	NO,	NZ,	OM,	
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ΑU	2003	2275	20		B9		2008	0626										
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CN	1655	787			A		2005	0817		CN 2	003-	8123	76		2	0030	527	

JP	2005531594	T	20051020	JP	2004-508810		20030527
NZ	536415	A	20061027	NZ	2003-536415		20030527
CN	101444509	A	20090603	CN	2008-10176372		20030527
US	20050239824	A1	20051027	US	2004-515275		20041122
US	7459464	B2	20081202				
MX	2004011859	A	20050331	MX	2004-11859		20041129
PRIORITY	APPLN. INFO.:			DK	2002-832	A	20020530
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OTHER SOURCE(S): MARPAT 140:13076

ΙT 195875-87-7

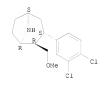
> RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

(triple monoamine reuptake inhibitors for treatment of chronic pain)

RN 195875-87-7 HCAPLUS

CN 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R, 2R, 3S, 5S) - (CA INDEX NAME)

Absolute stereochemistry.



OS.CITING REF COUNT: THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD (1 CITINGS)

REFERENCE COUNT: 3

THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD, ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 4 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:434366 HCAPLUS

DOCUMENT NUMBER: 139:963

Tropane derivatives having dopamine reuptake inhibitor TITLE: activity for the treatment of ischemic diseases

Scheel-Kruger, Jorgen; Ronn, Lars Christian B.

PATENT ASSIGNEE(S): Neurosearch A/S, Den. SOURCE: PCT Int. Appl., 42 pp.

CODEN: PIXXD2 DOCUMENT TYPE: Patent

LANGUAGE: English FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

INVENTOR(S):

PA:	PATENT NO.					D	DATE			APPLICATION NO.						DATE		
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WO	WO 2003045388					A1 20030605				WO 2	002-	DK79	6		20021128			
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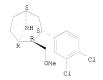
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8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-,

Absolute stereochemistry.

(1R, 2R, 3S, 5S) - (CA INDEX NAME)

CN



OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD (5 CITINGS)

REFERENCE COUNT: 10 THERE ARE 10 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L12 ANSWER 5 OF 5 HCAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1997:579720 HCAPLUS DOCUMENT NUMBER: 127:262895

ORIGINAL REFERENCE NO.: 127:51357a,51360a

TITLE: Preparation of tropane derivatives as inhibitors of monoamine neurotransmitter re-untake in the central

nervous system.

INVENTOR(S): Scheel-Kruger, Jorgen; Moldt, Peter; Watjen, Frank
PATENT ASSIGNEE(S): Neurosearch A/S, Den.; Scheel-Kruger, Jorgen; Moldt,

Peter; Watjen, Frank

SOURCE: PCT Int. Appl., 44 pp. CODEN: PIXXD2

DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9730997	A1	19970828	WO 1997-EP850	19970221

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T 20010715 AT 1997-903355

B6 20010806 SK 1998-929

A1 20010905 EP 2001-108256
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SG 99853 A1 20031127 SG 1999-3902
EE 4751 B1 20061215 EE 1998-254
US 6288079 B1 20010911 US 1998-101524
BG 63945 B1 20030731 BG 1998-102637
NO 9803877 A 19980821 NO 1998-3877
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HK 1018957 A1 20020927 HK 1999-104015
US 20010018444 A1 20010830 US 2001-814413
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W 19970221
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OTHER SOURCE(S): MARPAT 127:262895

IT 195875-87-7P

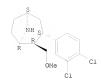
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of tropane derivs. as inhibitors of monoamine neurotransmitter re-uptake in central nervous system)

RN 195875-87-7 HCAPLUS

N 8-Azabicyclo[3.2.1]octane, 3-(3,4-dichlorophenyl)-2-(methoxymethyl)-, (1R,2R,3S,5S)- (CA INDEX NAME)

Absolute stereochemistry.



OS.CITING REF COUNT:

27 THERE ARE 27 CAPLUS RECORDS THAT CITE THIS

RECORD (32 CITINGS)

REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 54.90

94.89

FULL ESTIMATED COST

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